Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE011416\
Data File : BE091363.D
Acq On : 14 Jan 2016 23:00
Operator : UM/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE011416

Quant Time: Jan 15 15:15:23 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE011416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jan 14 23:29:02 2016
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	80	0.00
2	S	2-Fluorophenol	1.000	1.309	-30.9#	108	0.00
3	S	Phenol-d6	1.000	1.275	-27.5#	130	0.00
4	I	Naphthalene-d8	5.000	5.000	0.0	84	0.00
5	S	Nitrobenzene-d5	1.000	1.046	-4.6	88	0.00
6		Naphthalene	1.000	1.057	-5.7	87	0.00
7		2-Methylnaphthalene	1.000	0.955	4.5	100	0.00
8	I	Acenaphthene-d10	5.000	5.000	0.0	89	0.00
9	S	2,4,6-Tribromophenol	1.000	0.982	1.8	103	0.00
10	S	2-Fluorobiphenyl	1.000	1.078	-7.8	102	0.00
11		Acenaphthylene	1.000	0.961	3.9	103	0.00
12		Acenaphthene	1.000	1.031	-3.1	93	0.00
13		Fluorene	1.000	1.025	-2.5	104	0.00
14	I	Phenanthrene-d10	5.000	5.000	0.0	91	0.00
15		Phenanthrene	1.000	1.014	-1.4	90	0.00
16		Anthracene	1.000	0.959	4.1	99	0.00
17		Fluoranthene	1.000	0.812	18.8	69	0.00
18	I	Chrysene-d12	5.000	5.000	0.0	82	0.00
19		Pyrene	1.000	0.906	9.4	91	0.00
20	S	Terphenyl-d14	1.000	0.962	3.8	96	0.00
21		Benzo(a)anthracene	1.000	0.901	9.9	93	0.00
22		Chrysene	1.000	1.062	-6.2	87	0.00
23		Indeno(1,2,3-cd)pyrene	1.000	0.930	7.0	89	0.00
24	I	Perylene-d12	5.000	5.000	0.0	75	0.00
25		Benzo(b)fluoranthene	1.000	0.994	0.6	96	0.00
26		Benzo(k)fluoranthene	1.000	0.977	2.3	84	0.00
27	C	Benzo(a)pyrene	1.000	0.860	14.0	68	0.00
28		Dibenzo(a,h)anthracene	1.000	0.967	3.3	96	0.00
29		Benzo(g,h,i)perylene	1.000	0.945	5.5	84	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 0